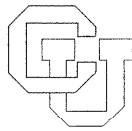


**Generalized Physical Networks for Automated Model Building \***

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**CU-CS-878-99**



**University of Colorado at Boulder  
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\* Supported by NSF NYI #CCR-9357740, ONR #N00014-96-1-0720, and a Packard Fellowship in Science and Engineering from the David and Lucille Packard Foundation.

Report Documentation Page			Form Approved OMB No. 0704-0188		
Public reporting burden for the collection of information is estimated to average 1 hour per response, including the time for reviewing instructions, searching existing data sources, gathering and maintaining the data needed, and completing and reviewing the collection of information. Send comments regarding this burden estimate or any other aspect of this collection of information, including suggestions for reducing this burden, to Washington Headquarters Services, Directorate for Information Operations and Reports, 1215 Jefferson Davis Highway, Suite 1204, Arlington VA 22202-4302. Respondents should be aware that notwithstanding any other provision of law, no person shall be subject to a penalty for failing to comply with a collection of information if it does not display a currently valid OMB control number.					
1. REPORT DATE <b>1999</b>		2. REPORT TYPE		3. DATES COVERED <b>00-00-1999 to 00-00-1999</b>	
4. TITLE AND SUBTITLE <b>Generalized Physical Networks for Automated Model Building</b>				5a. CONTRACT NUMBER	
				5b. GRANT NUMBER	
				5c. PROGRAM ELEMENT NUMBER	
6. AUTHOR(S)				5d. PROJECT NUMBER	
				5e. TASK NUMBER	
				5f. WORK UNIT NUMBER	
7. PERFORMING ORGANIZATION NAME(S) AND ADDRESS(ES) <b>Department of Computer Science, University of Colorado, Boulder, CO, 80309</b>				8. PERFORMING ORGANIZATION REPORT NUMBER	
9. SPONSORING/MONITORING AGENCY NAME(S) AND ADDRESS(ES)				10. SPONSOR/MONITOR'S ACRONYM(S)	
				11. SPONSOR/MONITOR'S REPORT NUMBER(S)	
12. DISTRIBUTION/AVAILABILITY STATEMENT <b>Approved for public release; distribution unlimited</b>					
13. SUPPLEMENTARY NOTES					
14. ABSTRACT <b>see report</b>					
15. SUBJECT TERMS					
16. SECURITY CLASSIFICATION OF:			17. LIMITATION OF ABSTRACT	18. NUMBER OF PAGES <b>10</b>	19a. NAME OF RESPONSIBLE PERSON
a. REPORT <b>unclassified</b>	b. ABSTRACT <b>unclassified</b>	c. THIS PAGE <b>unclassified</b>			

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ACKNOWLEDGMENTS SECTION.



# Generalized Physical Networks for Automated Model Building

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Technical Report #CU-CS-878-99

## Abstract

We present a new knowledge representation and reasoning framework for modeling nonlinear dynamical systems. The goals of this framework are to smoothly incorporate varying levels of domain knowledge and to tailor the reasoning methods — and hence the search space — accordingly. Our solution exploits generalized physical networks (GPN), a meta-level representation of idealized two-terminal elements, together with a hierarchy of qualitative and quantitative analysis tools, to produce a dynamic modeling domain whose complexity naturally adapts to the amount of available information about the target system.

## 1 Introduction

*System identification* (SID) is the process of identifying a dynamic model of an unknown system. The topic of this paper is a new knowledge representation and reasoning (KRR) framework that makes it possible to automate this kind of analysis. The challenges involved are significant. Applications in different fields of science and engineering demand different kinds of models and modeling techniques. Large aerospace structures, for example, usually have a variety of inputs (wind speed, engine thrust, and turbulence) and outputs (velocity, stress and strain indicators from wing sensors, etc.) and are described by partial differential equations; a simple electrical circuit, on the other hand, may have a single current source as an input and a voltmeter as an output, and can be described by an ordinary differential equation (ODE). Any representation designed for reasoning about models of such systems has to be both flexible enough to handle various degrees of uncertainty and complexity, and yet powerful enough to deal with situations in which the input signal may or may not be controllable.

System identification entails two steps, as shown in Fig. 1: *structural identification*, wherein one ascer-

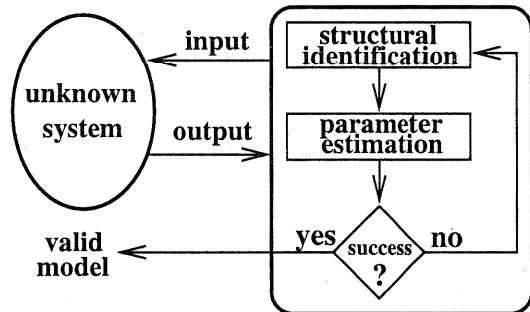


Figure 1: The system identification process

tains the general form of the model (e.g., the ODE  $a_1\ddot{\theta} + a_2\sin\theta = 0$  for a simple pendulum), and then *parameter estimation*, in which one finds specific parameter values for the unknown coefficients that fit that model to the observed data (e.g.  $a_1 = 1.0$ ,  $a_2 = -98.0$ ). Both steps depend upon *input-output analysis*[Casdagli, 1992], wherein the relationship between drive and response is used to infer useful information about internal system dynamics. For nonlinear systems, parameter estimation is difficult and structural identification is even harder; AI techniques can be used to automate the former[Bradley *et al.*, 1998], but the latter has, until now, remained the purview of human experts.

A central problem in any automated modeling task is that model complexity — and hence the size of the search space — is exponential in the number of model fragments unless severe restrictions are placed on the model-building process. Ideally, one would like to build *black-box* models using general reasoning techniques that applied to any system and did not require any *domain knowledge* about the system under examination. The combinatorics of the generate phase make this an unrealistic paradigm. A useful alternative is *gray-box* modeling, where the “box” is not completely opaque and information about its internals can be used to prune the search space down to a reasonable size. Some automated modeling tools, for example, maintain a library of typical components for a particular domain and build models by trying out various combinations of those primitives[Capelo

\*Supported by NSF NYI #CCR-9357740, ONR #N00014-96-1-0720, and a Packard Fellowship in Science and Engineering from the David and Lucile Packard Foundation.

*et al.*, 1998]. This approach is useful in relatively simple domains where the relationship between idealized modeling components and actual physical components is well known. Most AI modeling work has taken a *clear-box* modeling approach, in which one knows almost everything about what one is trying to model. The goal of the work described in this paper is to develop a KRR framework that supports automated modeling in a range of gray shades that is useful to practicing engineers.

The key to making gray-box modeling of nonlinear dynamical systems practical is a flexible knowledge representation scheme that adapts to the domain generality. Domain-dependent knowledge can reduce search-space size, but its applicability is fundamentally limited; the challenge in balancing these influences is to be able to determine, at every point in the reasoning procedure, what knowledge is (1) applicable and (2) useful. Knowledge that a particular input voltage controls an output frequency, for instance, allows one to reason more effectively about manipulating that voltage to learn more about the system.

The solution proposed here combines a representation that allows for different levels of domain knowledge, a set of reasoning techniques appropriate to each level, and a control strategy that invokes the right technique at the right time. In particular, we combine ideas from *generalized physical networks* [Sanford, 1965], a meta-level representation of idealized two-terminal elements, with traditional compositional model building [Falkenhainer and Forbus, 1991] and qualitative reasoning [Weld and de Kleer, 1990]. The intent is to bridge the gap between highly specific KRR frameworks that work well in a single, limited domain (e.g., a spring/dashpot vocabulary for modeling simple mechanical systems) and abstract frameworks that rely heavily upon general mathematical formalisms at the expense of having huge search spaces [Bradley and Stolle, 1996]. The generalized physical network (GPN) representation is an effective way to construct a description that is appropriate to a wide range of points on this spectrum, and thus it provides a useful bridge between general and specific modeling approaches.

## 2 Generalized Physical Networks

In the late 1950s and early 1960s, inspired by the realization that the principles underlying Newton’s third law and Kirchhoff’s current law were identical<sup>1</sup>, researchers began combining multi-port methods from a number of engineering fields into a generalized engineering domain with prototypical components [Paynter, 1961]. The basis of this *generalized physical networks* (GPN) paradigm is that the behavior of an ideal two-terminal element — the “component” — may be described by a mathematical relationship between two dependent variables: generalized flow and generalized effort, where  $flow(t) * effort(t) =$

$power(t)$ . This pair of variables manifests differently in each domain: ( $flow$ ,  $effort$ ) is ( $current$ ,  $voltage$ ) in an electrical domain and ( $force$ ,  $velocity$ ) in a mechanical domain<sup>2</sup>. One of the strengths of the GPN representation is that it brings out similarities between components and properties in different domains. Electrical resistors ( $v = iR$ ) and mechanical dampers ( $v = fB$ ) are analogous, as both dissipate energy. Similar relationships exist for generalized inertia, capacitance, flow, and effort source components, as shown in Table 1; see [Karnopp *et al.*, 1990] or [Sanford, 1965] for mechanical rotational, hydraulic, and thermal domains. Relationships also exist for pairs of mutually coupled two-terminal elements, such as electrical transformers or differential gears, in which a flow or effort in one element induces a flow or effort in the other. Components may be connected in the standard network-theoretic ways — parallel, serial, star, delta, etc. — and the network topology need not be planar.

Component	Electrical	Mechanical Translation
Resistor	$v = Ri$	$v = Bf$
Capacitor	$i = C \frac{dv}{dt}$	$f = M \frac{dv}{dt}$
Inertia	$v = L \frac{di}{dt}$	$v = K \frac{df}{dt}$
NL Resistor	$v = -R_A i^3$	$v = -R_A f^3$

Table 1: Example component representations.

Utilizing a small number of these primitive elements, GPNs can effectively model systems in a wide variety of domains; moreover, they make it very easy to incorporate varying amounts of information about those domains. A GPN-based modeling domain consists of a set of components and a set of connection primitives (“connectors”); a GPN model is a specific instance of connected components — e.g., a series connection of an inertia, a resistor, and a capacitor — represented by an incidence matrix. The same GPN model can be abstract and general or highly specific, depending on how much one knows about the domain. Both of the networks in Fig. 2, for example, can be modeled by a series inertia/resistor/capacitor GPN; knowledge that the system is electronic or mechanical would let one refine the model accordingly. The available domain knowledge, then, can be viewed as a lens that expands upon the internals of some GPN components, selectively sharpening the model *in appropriate and useful ways*.

There are a variety of ways to use generalized physical networks to help automate the structural identification phase of the SID process. One could, for example, create a library of all known components, enumerate all possible component combinations/configurations, and test each member of this succession until a valid model is

<sup>1</sup>Summation of {forces, currents} at a point is zero, respectively; both are manifestations of the conservation of energy.

<sup>2</sup>In bond graphs [Karnopp *et al.*, 1990], another generalized representation paradigm, velocity is a flow variable and force is an effort variable. The difference between GPNs and bond graphs is only a frame-of-reference shift.

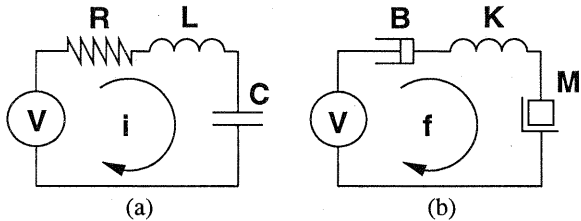


Figure 2: Two systems that are described by the same GPN model: (a) a series RLC circuit (b) a mass-spring-damper system.  $V$  is a voltage source in (a) and is a velocity source in (b).

found. This method is obviously impractical, as simple enumeration creates an exponential search space — a severe problem if the component library is large, as must be the case if one is attempting to model nonlinear systems<sup>3</sup>. A more-intelligent idea is to use a hierarchy of domain-dependent and -independent knowledge to direct the search, as described in the following section.

### 3 Reasoning with GPNs

The GPN representation, as described in the previous section, is an effective basis for a dynamic modeling domain whose complexity naturally rises and falls according to the available information about the target system. A general domain — e.g.,  $\{all\ dynamical\ systems\}$  — has a complex search space; a specific domain like  $\{conservative\ mechanical\ systems\}$  has a much smaller one. The challenge in reasoning about GPN models is to tailor the reasoning to the knowledge level in such a way as to prune the search space to the minimum. Organizing domains into a hierarchy of generality — e.g.,  $\{dynamical\ systems, electromechanical\ systems, mechanical\ systems, conservative\ mechanical\ systems, \dots\}$  — is not enough; what is needed is a hierarchical set of analysis tools, as well as a means for assessing the situation and choosing the appropriate tool.

Naively speaking, the goal is to start with a “plausible” model and then move through the space in a “reasonable” fashion. The key to doing so is to classify model and system behavior at an appropriate qualitative level. The order of an ODE, for instance, contains some useful high-level information about its behavior, and that kind of simple symbolic information can be used to remove huge branches from the search space. Sometimes, however, pruning a single leaf off the search tree requires expensive operations like point-by-point comparisons of ODE solutions to sensor data sets. Orchestrating this process efficiently is a difficult problem; its solution requires powerful machinery like theorem-provers and declarative meta-level control [Stolle and Bradley, 1998]. The generality of any behavior classification tool is also

<sup>3</sup>Nonlinear terms are somewhat idiosyncratic, and each must be supplied as a separate library entry. This issue has not arisen in previous work on GPNs because their use has been generally confined to linear systems.

affected by the domain of the target system; phase-portrait and time-series analysis apply to all ODEs, for instance, but transient analysis designed to detect “irrecoverable viscous deformation” is only meaningful for viscoelastic systems.

The KRR framework described in this paper is designed to exploit domain knowledge in order to navigate effectively through the search space of ODE models. The basic idea is that different reasoning techniques are appropriate to different domains, and the key to the solution is that the GPN component type and domain knowledge dictate which one to use. Table 2 shows part of the hierarchy of analysis tools that we have implemented. The top category is the broadest and most general. Lin-

System Type	Analysis Tools
Nonlinear	Cell-to-cell-mapping Delay-coordinate embedding NL time series analysis
Linear	Linear system analysis
Restricted linear	Domain dependent

Table 2: Component type and domain knowledge dictate what analysis tools apply. Tools higher in this table are more general but their results are less detailed.

ear systems are a subset of nonlinear systems; restricted linear systems — which use distributed parameters to model systems in very specific physical domains<sup>4</sup> — are, in turn, subsets of linear systems. Analysis tools that apply to any of these groups obviously apply to its subsets as well. All of the nonlinear systems tools listed in Table 2 work with the phase-space representation. *Cell-to-cell mapping* [Hsu, 1987] is a geometric reasoning technique that classifies a phase portrait qualitatively using simple discretized heuristics. *Delay-coordinate embedding* [Sauer *et al.*, 1991] lets one infer the dimension and topology of the internal system dynamics from a time series measured by a *single* output sensor. Nonlinear time-series analysis is a blanket term for classification that follows the  $\{attractors, bifurcations, \dots\}$  ontology of nonlinear dynamics [Strogatz, 1994]. Linear systems analysis refers to the techniques taught to undergraduate engineers (step/frequency response, etc.). Analysis tools for restricted linear systems — e.g., creep testing — are highly domain-specific.

Reasoning about models proceeds in the obvious manner, given this hierarchy: if no domain knowledge about the target system is available (i.e., the true “black-box” situation), then models are constructed, analyzed, and compared to the target system using general reasoning techniques that apply to *all* ODEs. Constructing models in this highly general manner is difficult, as there is no domain knowledge to limit the search, but it is also uni-

<sup>4</sup>e.g. modeling electrical transmission-line loss through series connections of multiple RC circuits; see Section 5 for an example in the viscoelastic domain.

versally applicable. If the system is known to be linear, the extensive and powerful repertoire of linear analysis tools developed over the last several decades makes the task far less imposing. Among other things, system inputs (drive terms) in linear systems appear verbatim in the resulting system ODE, which makes input/output analysis much easier. In restricted linear domains, analysis tools are even more specific and powerful. In viscoelastics, for example, three qualitative properties of a “strain test” reduce the search space of possible models to linear.

Given all of this machinery, we build models as follows. First, a candidate GPN model is constructed using a basic subset of the components and connectors of a particular domain. This process is guided by the analysis tools in Table 2. The cohort of nonlinear tools, for example, is first applied to the sensor data to determine the dimension  $d$  of the dynamics; this fact allows the modeling tool to automatically disregard all models of order  $< d$ . We have currently implemented roughly a dozen other nonlinear analysis techniques that yield similar search-space reductions; if the system is linear, several dozen more tools apply as well. Knowledge that the target system is oscillating, for example, not only constrains the model to be of least second order, but also implies some constraints on its coefficients. If this reduced search space does not contain a model that matches the observed system behavior, the GPN modeling domain is dynamically expanded to include more esoteric components. For example, if all models in the  $\{\text{linear-inductor}, \text{linear-resistor}\}$  domain are rejected, then the domain might be expanded to include *linear-capacitors*. Once a successful GPN network is found, the final step in the model-building process is to convert it into ODE format, as described in the following section.

## 4 Converting GPNs to ODEs

The last stage in the system identification process is to convert the domain-independent GPN model into ODE form. This conversion step is fairly easy to automate. The powerful network-theoretic principles involved have been in the engineering vernacular for many decades, but we use them in a very different manner. Traditionally, one applies a tool like *modified nodal analysis*<sup>5</sup> to a known network with known component values in order to analyze or simulate its behavior. We use loop and node equations to convert a GPN network with unspecified component values into an ODE (also with unspecified component values). Finally, we use a nonlinear parameter estimator [Bradley *et al.*, 1998] to find component values that match that ODE model to the observed system behavior. In this section, we describe how this procedure works for the electronics domain; the process is effectively identical in other domains.

The GPN→ODE algorithm is similar to standard system analysis: create a set of loop or node equations, substitute physical component models into those equations,

and then simplify/change variables as necessary. This algorithm is significantly simpler than a comparable bond graph→ODE conversion algorithm, since a bond graph algorithm creates implicit causal orderings as it creates the set of ODEs.

### Algorithm:

1. Use generalized Kirchhoff’s current law —  $\Sigma\{\text{flows into a node}\} = 0$  — to identify  $l$  independent branch currents,  $l = b - (n - 1)$ , where  $b$  is the number of branches and  $n$  is the number of nodes in the network. Branch independence is guaranteed by choosing each successive loop such that it traverses one and only one branch that is not part of an existing loop. The branches of a depth-first search tree of the network form one such set.
2. Use generalized Kirchhoff’s voltage law —  $\Sigma\{\text{efforts around a loop}\} = 0$  — around each of the  $l$  closed loops identified in step 1, yielding  $l$  system equations.
3. Apply the correct effort/flow relationships (from Table 1) to every element in the  $l$  loops from step 2, creating a set of ODEs.
4. Use variable substitution and/or symbolic differentiation to remove integrating elements as necessary.

As an example, consider a GPN modeling domain that allows parallel and serial connections of linear resistors (**R**), inductors (**L**), and nonlinear active resistors (**NR**) that obey the constitutive relation  $v = -R_A i^3 + R_B i$ . Midway through the procedure, after the modeler has tried and rejected a variety of too-simple models, we might see the candidate model shown in Fig. 3. Since

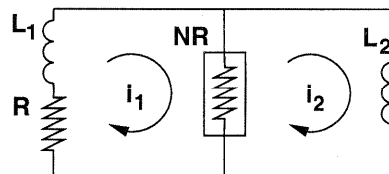


Figure 3: An electronics example.

$b = 4$  and  $n = 3$ , two loops are required. These are chosen automatically from the branches of a depth-first spanning tree.  $i_1$  and  $i_2$ , as shown, are one possibility;  $i_1$  and a loop through  $\{\mathbf{R}, \mathbf{L}_1, \mathbf{L}_2\}$  are another. Using generalized KVL around these two loops and substituting the given component models yields:

$$L_1 \frac{di_1}{dt} + R i_1 - R_A (i_1 - i_2)^3 + R_B (i_1 - i_2) = 0$$

$$L_1 \frac{di_2}{dt} - R_A (i_2 - i_1)^3 + R_B (i_2 - i_1) = 0$$

The last step is to pass this set of independent system equations in the two state variables  $\{i_1, i_2\}$  to the parameter estimator for comparison with observed behavior and determination of the unknown coefficients  $L_1, R,$

<sup>5</sup>which is based on node equations



etc. If the model fails this check, the modeling tool first tries to reconfigure it (e.g. moving  $L_1$  to be in series with  $NR$  instead of  $R$ ), and eventually gives up and expands the modeling domain (e.g. including linear capacitors). This *generate-and-test* process continues until a valid model is found.

This technique may initially appear to be a simulation/diagnostic tool for a system whose structure is completely known, but it is really much more abstract and adaptable. The KRR framework lets the modeling tool work with general GPN components as long as possible: up until the point when it converts the network into a set of ODEs. The relationships between component models and actual components of a physical system are a wholly separate data structure.

This approach has several important limitations and restrictions. Among other things, these models do not explicitly represent causality relationships — information that is very important for diagnostic tasks or automated explanation tools, but less so for system identification. The algorithm described above can also create extra system equations and variables; it may, for instance, construct a large resistive network when a single Thévenin equivalent would do. One can solve this problem using resistive network reduction [Mauss and Neumann, 1996]; we avoid it from the outset by building models in a sequential fashion and not adding like components in serial or parallel. Other meta-level representational schemes (such as bond graphs) solve these problems automatically, but GPNs are more expressive and more widely applicable, as well as easier to implement and use.

As is true of automated modeling in general, evaluating the results of this approach can be difficult because the question “How is this model better?” is hard to formalize. From an engineering standpoint, a successful model is one that matches observed behavior to within predefined specifications. PRET, the automated modeling tool in which these ideas are instantiated, is designed to be an engineer’s tool, so its judgement of what constitutes success or failure is exactly that. *Parsimony* is another desirable attribute in a model: one wishes to account for the observed behavior using as few — and as simple — ODEs terms as possible. Defining good metrics for model fragment simplicity is deceptively difficult, and it remains an open problem in automated modeling community. Finally, the speed with which PRET produces such a model is another important metric, particularly as we work with more-complex systems and search spaces. Ultimately, the best form of evaluation will consist of whether or not PRET’s models are useful for control system design — that is, whether the ODE that PRET constructs of a radio-controlled car can actually be used as the heart of a controller designed to direct that car to perform some prescribed action. We are in the process of evaluating models of real-world systems in several domains — ranging from robotics to hydrology — in this manner.

## 5 Related Work

Much of the pioneering work in the qualitative reasoning (QR) modeling community focuses on reasoning about pre-existing models: simulating them [Kuipers, 1986], simplifying and refining them [Weld, 1992], or keeping track of which model is appropriate in which regime [Ad-danki *et al.*, 1991]. QR model *construction* research has focused on building models from fragments [Falkenhainer and Forbus, 1991; Bobrow *et al.*, 1996; Williams and Nayak, 1996; Capelo *et al.*, 1998]. The work described in this paper uses some of the same techniques, but it has different goals and a different overall approach: it works with noisy, incomplete sensor data from real-world systems and attempts not to “discover” the underlying physics, but rather to find the simplest ODE that can account for the given observations.

Of all the QR modeling work cited in the previous paragraph, [Capelo *et al.*, 1998] is the most closely related to this paper. Capelo *et al.* build ODE models of linear viscoelastic systems by evaluating time series using qualitative reasoning techniques, and then use a parameter estimator to match the resulting model with a given system. This domain is limited — a “restricted linear system,” in our terminology — and fairly simple; it involves only two component types (linear springs and dashpots), connected in series and/or parallel. The specific nature of the domain naturally limits the search space, so a simple qualitative pre-processing step can be used to identify the system as one of only four different model types. The system described in this paper is much more general; it works on *all* linear and *nonlinear* lumped-parameter continuous-time ODEs and uses *dynamic* model generation to handle arbitrary devices and connection topologies in multiple domains. (Indeed, one of the domains that we have implemented is *{viscoelastic systems}*.)

The representation most closely related to GPNs is the bond graph [Karnopp *et al.*, 1990]. In the QR literature, bond graphs have been used primarily for reasoning about causality [Top and Akkermans, 1991] and modeling hybrid systems [Mosterman and Biswas, 1997]. An important exception to this is Amsterdam’s work on automated model construction in multiple physical domains [Amsterdam, 1992], which uses bond graphs to model linear systems of order two or less. While bond graphs are a good alternative to generalized physical networks — especially if causality issues are a concern — converting them into ODE models is difficult, which makes them less useful for the kinds of complex nonlinear modeling tasks that we address in this paper.

## 6 Conclusions and Future Work

Generalized physical networks, coupled with a hierarchy of qualitative and quantitative reasoning tools that relate observed physical behavior and model form, provide the flexibility required for gray-box modeling of nonlinear dynamical systems. This type of reasoning, wherein the modeling tool has only partial knowledge of the in-

ternals of the target system, accurately reflects the abstraction levels and reasoning processes used effectively by human engineers during the modeling procedure. The KRR framework described in this paper allows automated modeling tools to reason effectively with varying levels of domain knowledge — about different domains and at different levels of abstraction in an individual domain. The GPN component representation adapts naturally to the information available about the target system, allowing for the creation of dynamic modeling domains. Detailed knowledge automatically reduces the search space of models and triggers powerful, specialized reasoning techniques that are appropriate for that situation; in the absence of such knowledge, the system falls back on broadly applicable domain-independent principles to navigate through an exponential search space.

Besides speeding model generation, this KRR framework streamlines reasoning about the relationship between input sources and output responses. We are currently working on automating this kind of reasoning, using GPNs to represent sensors and actuators explicitly with the goal of automatically performing experiments on unknown systems. We are also investigating model fragments and model complexity. Our domain libraries currently include linear components and some simple nonlinear ones; if the modeling task calls for a more-complicated device, we first query the user and then resort to power-series expansions. Once we have experimented with more nonlinear systems in more domains, we may augment our libraries, and it may be useful to organize them in hierarchies (e.g. one linear spring as a starting point, with a simple nonlinear spring as an alternative, and more-complicated devices further down the hierarchy).

**Acknowledgments:** Reinhard Stolle, Joe Iwanski, Apollo Hogan, and Brian LaMacchia also contributed code and/or ideas to this project, and the IJCAI reviewers' comments helped focus the content of this paper.

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